

The Magazine for ESRI Software Users

July-September 2007 Vol. 10 No. 3

Optimizing a Sampling Network Automating the Use of Geostatistical Tools for Lake Tahoe Area Study

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Editor's note: This article describes how powerful analysis tools in the ArcGIS Geostatistical Analyst 9.2 extension were applied in a study of air quality degradation around Lake Tahoe, a resort destination located on the California/Nevada border. As part of the study, a model to optimize the monitoring network by locating additional monitoring stations was built using ModelBuilder. An accompanying article, "Making Effective Use of Geostatistics," introduces this class of statistics.

About the Study

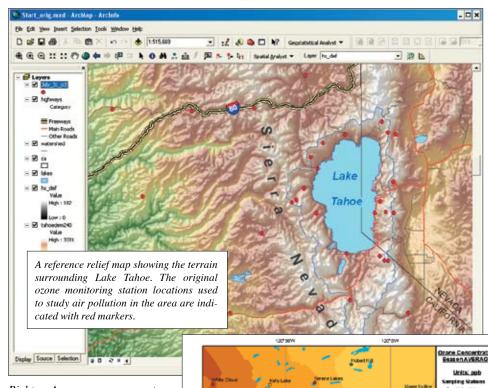
The transparency and purity of Lake Tahoe's water has been deteriorating since the 1950s, partially due to increased deposition of nitrogenous air pollutants. Forests in the Lake Tahoe watershed have also suffered from stresses such as drought, overstocking, and elevated concentrations of phytotoxic air pollutants, mainly ozone.

Ozone, one of the most damaging air pollutants, has strong toxic effects on human health and vegetation and is most indicative of photochemical smog.

One of the main questions for scientists and forest managers in this area is whether air pollution (specifically ozone) is generated locally or is migrating with the prevailing westerly winds from California's Central Valley, an area known for high levels of air pollution. A study to determine the origin of ozone found in the vicinity of Lake Tahoe was undertaken by the Forest Service and ESRI. Initially, ambient ozone concentrations were measured using a network of 31 sampling stations established by the United States Department of Agriculture (USDA) Forest Service Pacific Southwest Research Station scientists from the Riverside Fire Laboratory in Riverside, California.

Looking at the Original Network

Knowledge of the central Sierra Nevada Mountains, both the general patterns of ozone distribution and the westerly wind pattern in this area, led to the positioning of several monitoring points on the western slopes of these mountains. Most monitoring stations were located inside the Lake Tahoe Basin.



Right: An ozone concentration prediction map showing the west-east trend in ozone concentration.

Three elevation transects were set to measure ozone concentration at different altitudes to learn if these concentrations could be correlated with elevation.

Prediction maps of ozone concentration were generated using Geostatistical Analyst. Because no strong correlation was de-

tected between ozone and elevation, cokriging could not be applied. Most maps of ozone concentration showed a noticeable trend. The main range of the Sierra Nevada Mountains apparently blocks the transport of ozone from the Central Valley, located to the west of the study area. This helps explain why no correlation between ozone concentration and elevation was detected. Monitoring stations located at similar elevations

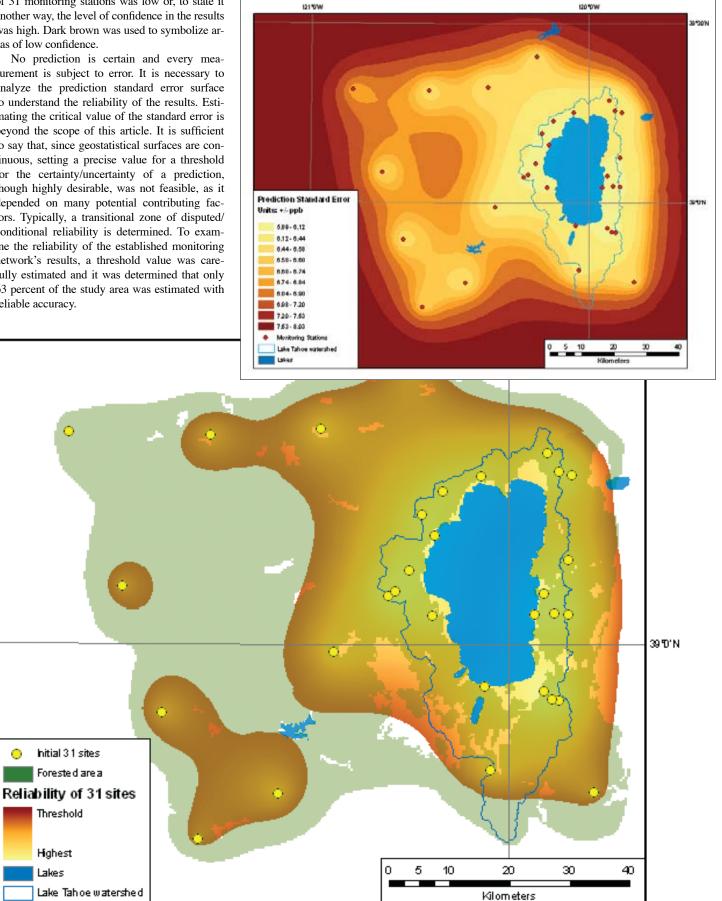
but on opposite sides of the main range reported significantly different ozone values.

Looking at Error

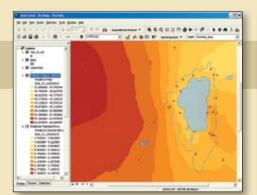
A map of prediction standard error was created using the same kriging method and parameters that were used to generate ozone prediction maps. The bright yellow colors indicated areas where the prediction standard error for the existing network

of 31 monitoring stations was low or, to state it another way, the level of confidence in the results was high. Dark brown was used to symbolize areas of low confidence.

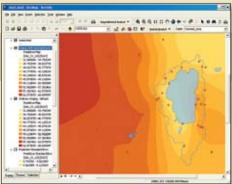
No prediction is certain and every measurement is subject to error. It is necessary to analyze the prediction standard error surface to understand the reliability of the results. Estimating the critical value of the standard error is beyond the scope of this article. It is sufficient to say that, since geostatistical surfaces are continuous, setting a precise value for a threshold for the certainty/uncertainty of a prediction, though highly desirable, was not feasible, as it depended on many potential contributing factors. Typically, a transitional zone of disputed/ conditional reliability is determined. To examine the reliability of the established monitoring network's results, a threshold value was carefully estimated and it was determined that only 63 percent of the study area was estimated with reliable accuracy.



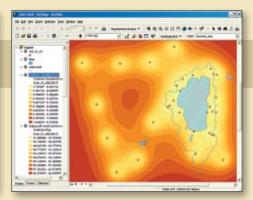
Above right: The prediction standard error map generated with ordinary kriging shows surfaces of prediction uncertainty classified into 10 categories with initial monitoring stations labeled. Bottom: The study area with the forested area surrounding Lake Tahoe. The initial 31-station network did not adequately cover this area.



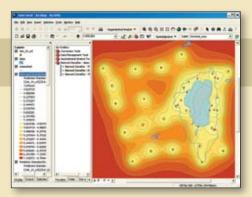
This prediction map of ozone concentrations was made using ordinary kriging and software default values.



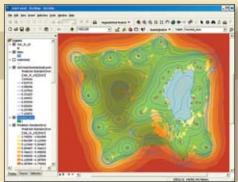
To create a more reliable surface, trend and error modeling were added.



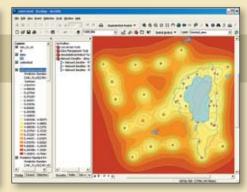
After evaluating the reliability of the surface, it was clear that more sampling points were needed.



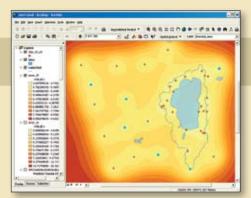
The increase in the yellow areas indicates that adding six monitoring stations has made the network more reliable.



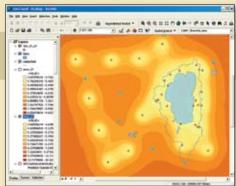
The output geostatistical surfaces of prediction standard error generated by the model were converted into rasters with the same color schema and displayed with the forest area.



The surface of prediction standard error based on 37 sampling points was considerably brighter, signifying that the prediction of ozone concentration based on the enhanced network would be more reliable.



The locations of the additional stations are indicated in cyan.



The prediction standard error surface based on a network of 37 points is much brighter and more reliable.

Making the Network More Reliable

Because the reliability of geostatistical analysis depends on having a sufficient number of appropriately distributed sampling stations, monitoring activities commonly encounter problems caused by networks of sampling points that are not sufficiently dense.

GIS can be applied to optimize a monitoring network. In this case, it was used to improve the reliability of the ozone concentration predictions. Models of ozone concentration and models of prediction standard error of ozone concentration were generated to determine locations where new sampling stations were most needed. These stations could be added until the surface of prediction standard error for the entire study area was below a given threshold or until the project budget was exhausted—whichever came first.

This article provides an overview of how Geostatistical Analyst was applied to that end rather than a detailed description of each step taken. The initial network of 31 points was improved using a model that was created in ModelBuilder. This Automated Network Densifier model incorporated new options in Geostatistical Analyst that were introduced with ArcGIS Desktop 9.2.

Two ozone prediction maps were generated and compared. The methods and parameters used to generate the one with higher accuracy were applied to create maps of prediction standard error. Several supplemental points were sequentially added to the network at the locations showing the lowest reliability. To graphically demonstrate the increasing trustworthiness of the predictions, based on the growing number of monitoring sites, maps of the variability of the prediction standard error used the same color symbology.

Using the Default Parameters

The study used ozone concentrations obtained for the month of July because during that month ozone concentrations are usually highly elevated and have greater potential for both harmful health effects and damage to forest vegetation.

Using the Geostatistical Wizard, a tool in Geostatistical Analyst that leads users through the process of creating a statistically valid surface, a prediction map was created for the ozone concentration. It was generated by applying the default options for ordinary kriging. The resulting map indicated that the spatial distribution of ozone in July over the study area had a west–east trend (i.e., a high concentration in the west, a low concentration in the east, and continuously changing intermediate concentrations in between).

While the map generated using default values was acceptable, it is always more desirable to produce distribution maps of a given phenomenon as accurately as possible. Geostatistical Analyst has many methods and options. Expertise and

experience are required for optimally generating a more reliable geostatistical surface. It is necessary to customize methods and parameters for each dataset to create a surface more reliable than the one generated by default values.

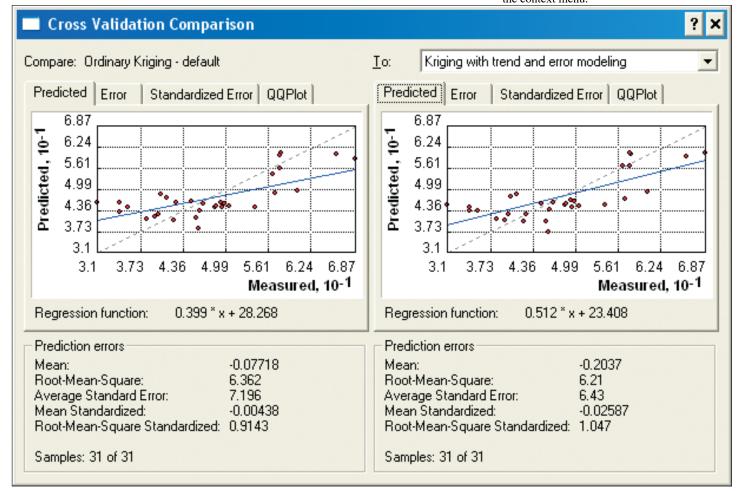
Modifying Default Parameters

For the ozone dataset, the following kriging parameters were applied.

- Const option of the order of trend removal was set to 67 percent Global.
- The Smooth option was used to make the appearance of the output map look smoother.
- The Nugget effect value was reduced to 15.
- The Major Range was set to 40,000.
- The Error Modeling slider for measurement error was set to 50 percent.

The new layer generated from these parameters was named Kriging with trend and error modeling. The resulting surface was smoother, yet more detailed. It showed more features of ozone distribution in addition to those generated using only the default parameters.

Despite noticeable differences, the Kriging with trend and error modeling surface visually resembled the previous one. Which surface was more reliable? Geostatistical Analyst has special tools that help the user select the best surface. These tools are accessed by right-clicking the newly generated geostatistical surface in the table of contents and selecting Compare from the context menu.



Performing Diagnostics

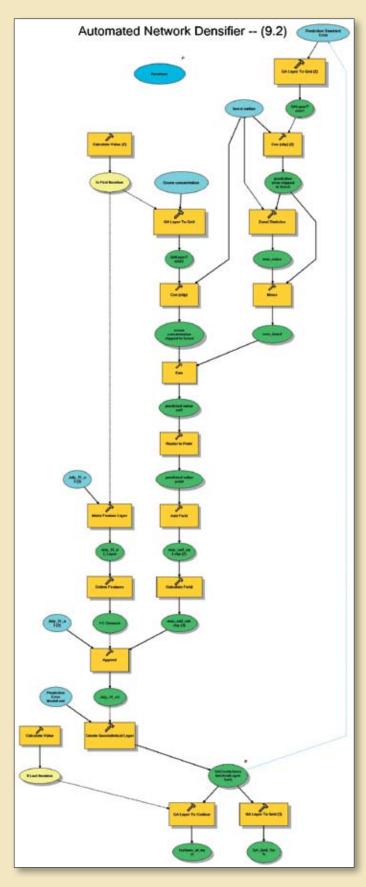
The Cross Validation Comparison dialog box compares five prediction error parameters. All error parameters, with the exception of the Root-Mean-Square Standardized, should be as close to 0 as possible for the most accurate output. For the Root-Mean-Square Standardized parameter, the result should be close to 1. In this example, four out of five of the error parameters indicated that the surface created by applying both trend and error modeling was more accurate. The most critical indicator of prediction accuracy, Root-Mean-Square Standardized, also indicated that the surface created using trend and error modeling was the winner. Obviously, additional prediction standard error surfaces could be generated using different sets of parameters.

After selecting the most accurate surface, the spatial distribution of the reliability of that layer (or in other words, the levels of uncertainty in the generated surface of ozone concentration) can be determined by creating a prediction standard error map. This map was generated using exactly the same parameters that were used for the latest prediction map. From the Method Summary interface, the methods and parameters that were used to generate the most accurate geostatistical surface were saved to an XML file for use later in the process. The same symbology was used-bright yellow color for the areas of highest reliability and dark brown for areas of lowest reliability. This prediction map showed that areas at the map edges and in the central part of the study area had low prediction reliability.

To improve the reliability of the ozone distribution surface and determine whether there was a significant ozone-generating source in the vicinity of Lake Tahoe, more sampling points were needed. The project's budget allowed for six additional measurement stations for the next season. Locations of new monitoring points were chosen to improve the overall reliability of the geostatistical interpolation by sampling at the locations within the study area where reliability was the lowest. In addition, all supplemental points had to be located within the forested portion of the study area.

Locations for the new points could be selected manually in ArcMap based on the criteria previously stipulated. Alternatively, locations could be selected using an automated method—a model. Part of the ArcGIS geoprocessing framework, ModelBuilder provides a graphic environment for creating, running, and saving models. Introducing a model would reduce subjectivity, make the selection of the prospective locations reproducible, and make the rules transparent.

The Automated Network Densifier model created for this project generated an enhanced monitoring network by adding supplemental points at the locations where they were most needed to reduce the overall prediction uncertainty. The prediction standard error geostatistical surface, the input data for the model, was



The Automated Network Densifier Model

The model converts the two geostatistical surfaces to the grid raster format and clips both to the geographic extent of the study area. The maximum value on the standard error of prediction grid is found, and that cell is converted into a point shapefile with a single feature and the value of the ozone concentration as its attribute. This is the optimal point for adding a station.

In the next iteration, this feature is appended to the original sampling network shapefile and a new geostatistical layer of prediction standard error is generated from the original 31 stations plus the new stations. The XML file containing the initial prediction standard error surface parameters is used to generate each *iteration. The process* is repeated to identify as many network sampling locations as desired.

In addition to five input datasets and 18 utilized functions, the model consists of two preconditions. First, since the ozone concentration geostatistical layer is not changed during the workflow, it is converted into a grid only during the first iteration. Second, only the last output grid of prediction standard error is converted into isolines of equal values of prediction standard error (contours).

updated at each iteration as the model appended one additional sampling point to the current monitoring network. The model could be run once to indicate where the most crucial missing point was located. It could also be run for a specified number of iterations to generate as many additional points as the project's budget allowed or until a variable was no longer equal to a predetermined condition. For example, it could be run until the maximum standard error of prediction for the study area was less than the largest acceptable potential error of prediction. The new points were sequentially placed at the location of the largest current potential standard error of prediction.

The Automated Network Densifier model takes five input data layers:

- The established monitoring network of 31 sites (as a shapefile)
- The geostatistical surface of ozone concentration that is based on the 31 original ozone monitoring sites
- The geostatistical surface of prediction standard error that is based on the original 31 monitoring sites
- The XML file containing the methods and parameters used to create the prediction standard error geostatistical layer originally selected
- The forested area at the vicinity of Lake Tahoe that constitutes the study area (as a grid file)

This model does not account for proximity to roads or access restrictions because adequate data for these factors was not available. Assuming the project budget allows for several monitoring stations to supplement the initial network, six iterations of the model were run and the locations for six new sampling stations were determined. The geostatistical surface of the standard error of prediction resulting from the sixth iteration was displayed together with the relevant vector version of the isolines of prediction standard error.

The points added to the network can be more easily seen by turning off the newly created layers in the table of contents and looking for the discrepancies between the geostatistical layer and the isolines of prediction trust. Because the supplemental points had to be located within the forested area, displaying the forested area grid as a semitransparent green polygon made it easier to understand why these locations were chosen.

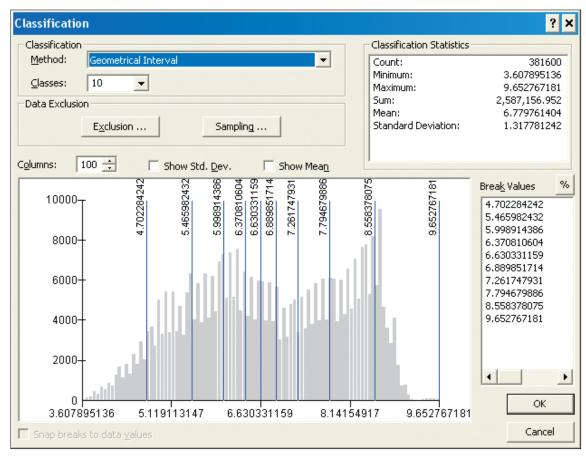
As expected, increasing the density of the monitoring network decreased the standard error of prediction for the entire study area. To better measure the increase in reliability caused by adding supplemental stations, the output geostatistical surfaces of prediction standard error generated by the model were converted into rasters with the same color schema.

In ArcGIS 9.2, users can now apply the Geometrical Interval classification to both rasters and geostatistical surfaces simply by right-clicking on the layer, choosing Properties > Symbology, and using the Classification option to change the classification to Geometric Interval. This method was applied to the output grid from the

final iteration of the Automated Network Densifier model with the number of classes set to 10 and a yellow to dark red color ramp.

Comparing the prediction standard error surface generated based on the initial 31 stations with the surface that used all 37 stations illustrated how the level of certainty of the prediction improved when the final grid was rendered using the same color symbology. The prediction standard error based on the original number of stations was displayed with the additional stations. The surface of prediction standard error based on 37 sampling points was considerably brighter, signifying that the prediction of ozone concentration based on the enhanced network would be more reliable. Without going into numerical details, the network of 37 stations can provide enough sampling data to significantly improve the trustworthiness of prediction over the entire study area. Whether the six additional stations for this network were sufficient to meet a minimum acceptable threshold of reliability is beyond the scope of this article.

The final result seems to confirm that the Automated Network Densifier model can improve the network design process during the second stage of sequential sampling. The proposed method appends the supplemental sites in a reasonable manner. It adds new points where they are most effective in enhancing reliability. With the model, as each new point is added, the improvement in the network can be observed.



Classification using Geometrical Interval is now available.

About the Authors

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Making Effective Use of Geostatistics

Applying GIS tools for studying the geographic distribution of regionalized variables

By Witold Fraczek, ESRI Application Prototype Lab, and Andrzej Bytnerowicz, USDA Forest Service

Geostatistics is a branch of science that applies statistical methods to spatial interpolation. Although geostatistics was developed independently of GIS, it has become an integral part of GIS. Without a computer and GIS mapping ability, it wouldn't be known outside a small group of geostatistical gurus. Just as one does not have to be a GIS expert to use GIS, one doesn't need to be a geostatistician to make effective use of geostatistics. Meteorologists, soil scientists, geologists, oceanographers, foresters, and other scientists can benefit from using appropriate geostatistical methods.

The functionality of geostatistics is applicable when the studied phenomena are regionalized variables that fall between random and deterministic variables. The geographic distribution of regionalized variables cannot be mathematically described as deterministic; yet the distribution of intensity of those phenomena is not random. Most of the natural phenomena that take place in the atmosphere, seawater, or soil meet the criteria of this category. The distribution of air temperature, the salinity of an ocean, soil moisture, or ore deposit concentrations in a geologic layer are examples of regionalized variables. Even though they don't represent truly natural phenomena, crop yield prediction and air pollution might also be subjects for geostatistical analysis.

It is not practical or possible to make exhaustive real-world observations so sampling is used for these analyses. The ultimate goal of sampling is to get a good representation of the phenomenon under study. Spatial sampling is an important consideration in environmental studies because sample configuration influences the reliability, effectiveness, and cost of a survey. Intensive sampling is expensive but gives a precise picture of spatial variability for a given phenomenon. However, sparse sampling is less expensive but may miss significant spatial features. Practical sampling constraints and the availability of existing information can enhance the development of a sampling scheme.

To ensure a high level of confidence in the results of any geostatistical interpolation, it is important to have a sufficient number of well-distributed sampling stations in the monitoring network. How many stations are sufficient and how can their distribution be optimized? GIS, and particularly the ArcGIS Geostatistical Analyst extension, can help answer this question.

One technique used to design an optimal sampling network for a regionalized variable, such as air pollution, is sequential sampling. Sequential sampling is based on extended knowledge of the area to be sampled and expertise in the factors controlling the distribution of a regionalized variable. Familiarity with the terrain and the phenomena should inform the initial choice of site for the sampling network. The results of this preliminary study are used to optimize the scheme by adding new sampling points both in areas having the lowest reliability and in possible hot spot areas (e.g., areas of maximum concentration, high variability, or uncertain measurements).

The kriging interpolator is considered the most sophisticated and accurate way to determine the intensity of a phenomenon at unmeasured locations. Kriging weights surrounding measured values are based not only on the distance between measured points and the prediction location but also on the overall spatial arrangement of the measured points. Except for generating an estimated prediction, kriging can provide a measure of an error, or uncertainty of the estimated surface. Since the estimation variances can be mapped, a confidence placed in the estimates can be calculated and their spatial distribution can be presented on a map to assist in the decision-making process. The prediction standard error maps show a distribution of a square root of a prediction variance, which is a variation associated with differences between the measured and calculated values. The prediction standard error quantifies an uncertainty of a prediction.



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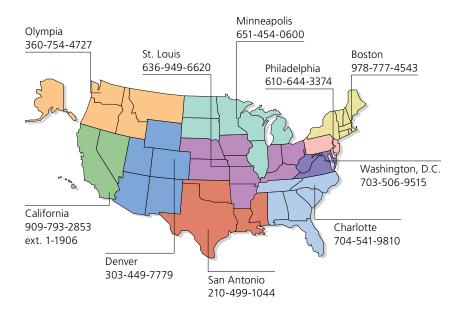
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